

USA2001/0081 US NP

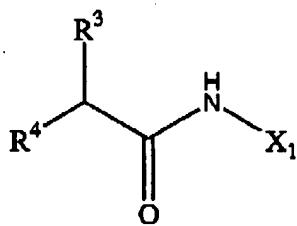
PATENT

Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently Amended) A compound of Formula I:



I

in which:

X¹ is -C(R¹)(R²)X²;

X² is cyano, -CHO, -C(R⁷)(R⁸)R⁵, -C(R⁷)(R⁸)CF₃, -C(R⁷)(R⁸)CF₂CF₂R⁹

-CH=CHS(O)₂R⁵, -C(R⁷)(R⁸)CF₂C(O)NR⁵R⁶, -C(R⁷)(R⁸)C(R⁷)(R⁸)NR⁵R⁶,

-C(R⁷)(R⁸)C(R⁷)(R⁸)OR⁵, -C(R⁷)(R⁸)CH₂OR⁵, -C(R⁷)(R⁸)CH₂N(R⁶)SO₂R⁵,

-C(R⁷)(R⁸)C(R⁷)(R⁸)N(R⁶)(CH₂)₂OR⁶, -C(R⁷)(R⁸)C(R⁷)(R⁸)N(R⁶)(CH₂)₂NR⁶ or

-C(R⁷)(R⁸)C(R⁷)(R⁸)R⁵; wherein R⁵ is (C₁₋₄)alkyl, (C₆₋₁₀)aryl(C₀₋₆)alkyl,

hetero(C₄₋₁₀)aryl(C₀₋₆)alkyl, (C₄₋₁₀)cycloalkyl(C₀₋₆)alkyl or hetero(C₄₋₁₀)cycloalkyl(C₀₋₆)alkyl

wherein hetero(C₄₋₁₀)aryl or hetero(C₄₋₁₀)cycloalkyl is pyran, thiopyran, pyrimidine, thiazole,

isothiazole, pyridine, furan, imidazole, isoxazole, oxadiazole, oxazole or triazole; R⁶ is hydrogen or (C₁₋₆)alkyl; R⁷ is hydrogen or (C₁₋₄)alkyl and R⁸ is hydroxy or R⁷ and R⁸ together form oxo;

R⁹ is hydrogen, halo, (C₁₋₄)alkyl, or (C₅₋₁₀)aryl(C₀₋₆)alkyl;

wherein within X¹ any cycloalkyl, is substituted or unsubstituted;

R¹ and R² are both fluoro; or

R¹ is hydrogen or (C₁₋₆)alkyl and R² is selected from the group consisting of hydrogen, (C₁₋₆)alkyl, cyano, -X⁴NR¹²R¹², -X⁴NR¹²C(O)R¹², -X⁴NR¹²C(O)OR¹², -X⁴NR¹²C(O)NR¹²R¹², -X⁴NR¹²C(NR¹²)NR¹²R¹², -X⁴OR¹³, -X⁴SR¹³, -X⁴C(O)OR¹², -X⁴C(O)R¹³, -X⁴OC(O)R¹³,

USA V2001/0081 US NP

PATENT

$-X^4C(O)NR^{12}R^{12}$, $-X^4S(O)_2NR^{12}R^{12}$, $-X^4NR^{12}S(O)_2R^{13}$, $-X^4P(O)(OR^{12})OR^{12}$,
 $-X^4OP(O)(OR^{12})OR^{12}$, $-X^4S(O)R^{14}$, $-X^4S(O)_2R^{14}$, $-R^{15}$, $-X^4OR^{15}$, $-X^4SR^{15}$, $-X^4S(O)R^{15}$,
 $-X^4S(O)_2R^{15}$, $-X^4C(O)R^{15}$, $-X^4C(O)OR^{15}$, $-X^4OC(O)R^{15}$, $-X^4NR^{15}R^{12}$, $-X^4NR^{12}C(O)R^{15}$,
 $-X^4NR^{12}C(O)OR^{15}$, $-X^4C(O)NR^{15}R^{12}$, $-X^4S(O)_2NR^{15}R^{12}$, $-X^4NR^{12}S(O)_2R^{15}$,
 $-X^4NR^{12}C(O)NR^{15}R^{12}$ and $-X^4NR^{12}C(NR^{12})NR^{15}R^{12}$, wherein X^4 is a bond or (C_{1-6}) alkylene, R^{12} at each occurrence independently is hydrogen or (C_{1-6}) alkyl, R^{13} is hydrogen, (C_{1-6}) alkyl or halo-substituted(C_{1-6})alkyl, R^{14} is (C_{1-6}) alkyl or halo-substituted(C_{1-6})alkyl, and R^{15} is (C_{3-10}) cycloalkyl(C_{0-6})alkyl, (C_{6-10}) aryl(C_{0-6})alkyl, or (C_{9-12}) bicycloaryl(C_{0-6})alkyl or ~~merpholinyl~~;
 $or R^1$ and R^2 taken together with the carbon atom to which both R^1 and R^2 are attached form (C_{3-8}) cycloalkylene; wherein R^2 , and said cycloalkylene may be substituted further with 1 to 3 radicals independently selected from (C_{1-6}) alkyl, cyano, halo, halo-substituted($C_{1-4})$ alkyl, nitro, $-X^4NR^{12}R^{12}$, $-X^4NR^{12}C(O)R^{12}$, $-X^4NR^{12}C(O)OR^{12}$, $-X^4NR^{12}C(O)NR^{12}R^{12}$, $-X^4NR^{12}C(NR^{12})NR^{12}R^{12}$, $-X^4OR^{13}$, $-X^4SR^{13}$, $-X^4C(O)OR^{12}$, $-X^4C(O)R^{13}$, $-X^4OC(O)R^{13}$, $-X^4C(O)NR^{12}R^{12}$, $-X^4S(O)_2NR^{12}R^{12}$, $-X^4NR^{12}S(O)_2R^{13}$, $-X^4P(O)(OR^{12})OR^{12}$, $-X^4OP(O)(OR^{12})OR^{12}$, $-X^4S(O)R^{14}$ and $-X^4S(O)_2R^{14}$, wherein X^4 , R^{12} , R^{13} and R^{14} are as defined above;

R^2 and R^4 are independently $-C(R^{16})(R^{17})X^7$, wherein R^{16} and R^{17} are hydrogen, (C_{1-6}) alkyl or fluoro, or R^{16} is hydrogen and R^{17} is hydroxy and X^7 is selected from $-X^4NR^{12}R^{12}$, $-X^4NR^{12}C(O)R^{12}$, $-X^4NR^{12}C(O)OR^{12}$, $-X^4NR^{12}C(O)NR^{12}R^{12}$, $-X^4NR^{12}C(NR^{12})NR^{12}R^{12}$, $-X^4OR^{13}$, $-X^4SR^{13}$, $-X^4C(O)OR^{13}$, $-X^4C(O)R^{13}$, $-X^4OC(O)R^{13}$, $-X^4C(O)NR^{12}R^{12}$, $-X^4S(O)_2NR^{12}R^{12}$, $-X^4NR^{12}S(O)_2R^{13}$, $-X^4P(O)(OR^{12})OR^{12}$, $-X^4OP(O)(OR^{12})OR^{12}$, $-X^4S(O)R^{14}$, $-X^4S(O)_2R^{14}$, $-R^{15}$, $-X^4OR^{15}$, $-X^4SR^{15}$, $-X^4S(O)R^{15}$, $-X^4S(O)_2R^{15}$, $-X^4C(O)R^{15}$, $-X^4C(O)OR^{15}$, $-X^4OC(O)R^{15}$, $-X^4NR^{15}R^{12}$, $-X^4NR^{12}C(O)R^{15}$, $-X^4NR^{12}C(O)OR^{15}$, $-X^4C(O)NR^{15}R^{12}$, $-X^4S(O)_2NR^{15}R^{12}$, $-X^4NR^{12}S(O)_2R^{15}$, $-X^4NR^{12}C(O)NR^{15}R^{12}$ and $-X^4NR^{12}C(NR^{12})NR^{15}R^{12}$, wherein X^4 , R^{12} , R^{13} , R^{14} and R^{15} are as defined above;

R^3 is $-C(R^{16})(R^{17})X^7$, wherein R^{16} and R^{17} are hydrogen, (C_{1-6}) alkyl or fluoro, or R^{16} is hydrogen and R^{17} is hydroxyl; and X^7 is selected from $-X^4NR^{12}R^{12}$, $-X^4NR^{12}C(O)R^{12}$, $-X^4NR^{12}C(O)OR^{12}$, $-X^4NR^{12}C(O)NR^{12}R^{12}$, $-X^4NR^{12}C(NR^{12})NR^{12}R^{12}$, $-X^4OR^{13}$, $-X^4SR^{13}$, $-X^4C(O)OR^{12}$, $-X^4C(O)R^{13}$, $-X^4OC(O)R^{13}$, $-X^4C(O)NR^{12}R^{12}$, $-X^4S(O)_2NR^{12}R^{12}$, $-X^4NR^{12}S(O)_2R^{13}$, $-X^4P(O)(OR^{12})OR^{12}$, $-X^4OP(O)(OR^{12})OR^{12}$, $-X^4S(O)R^{14}$, $-X^4S(O)_2R^{14}$, $-R^{15}$

USA V2001/0081 US NP

PATENT

$-X^4OR^{15}$, $-X^4SR^{15}$, $-X^4S(O)R^{15}$, $-X^4S(O)_2R^{15}$, $-X^4C(O)R^{15}$, $-X^4C(O)OR^{15}$, $-X^4OC(O)R^{15}$,
 $-X^4NR^{15}R^{12}$, $-X^4NR^{12}C(O)R^{15}$, $-X^4NR^{12}C(O)OR^{15}$, $-X^4C(O)NR^{15}R^{12}$, $-X^4S(O)_2NR^{15}R^{12}$,
 $-X^4NR^{12}S(O)_2R^{15}$, $-X^4NR^{12}C(O)NR^{15}R^{12}$ and $-X^4NR^{12}C(NR^{12})NR^{15}R^{12}$, wherein X^4 , R^{12} , R^{13} ,
 R^{14} and R^{15} are as defined above;

R^4 is $-C(R^{16})(R^{17})X^7$, wherein R^{16} and R^{17} are hydrogen, (C_{1-6}) alkyl or fluoro, or R^{16} is
hydrogen and R^{17} is hydroxy and X^7 is selected from $-R^{15}$, $-X^4OR^{15}$, $-X^4SR^{15}$, $-X^4S(O)R^{15}$,
 $-X^4S(O)_2R^{15}$, $-X^4C(O)R^{15}$, $-X^4C(O)OR^{15}$, $-X^4OC(O)R^{15}$, $-X^4NR^{15}R^{12}$, $-X^4NR^{12}C(O)R^{15}$,
 $-X^4NR^{12}C(O)OR^{15}$, $-X^4C(O)NR^{15}R^{12}$, $-X^4S(O)_2NR^{15}R^{12}$, $-X^4NR^{12}S(O)_2R^{15}$,
 $-X^4NR^{12}C(O)NR^{15}R^{12}$ and $-X^4NR^{12}C(NR^{12})NR^{15}R^{12}$, wherein X^4 , R^{12} , R^{13} and R^{14} are as defined
above for R^3 and R^{15} is hetero(C_{3-10})cycloalkyl(C_{0-3})alkyl wherein said
hetero(C_{3-10})cycloalkyl(C_{0-3})alkyl is morpholinyl,

wherein within one of R^3 or R^4 any cycloalkyl, or aryl or is substituted or unsubstituted and wherein each of R^3 and R^4 is substituted further or is not further substituted, and provided that when X^2 is cyano and X^7 within one of R^3 or R^4 is $-X^4C(O)R^{13}$ or $-X^4C(O)R^{15}$, wherein X^4 is a bond, then X^7 within the other of R^3 or R^4 is limited to $-X^4SR^{15}$, $-X^4S(O)R^{15}$ and $-X^4S(O)_2R^{15}$, wherein R^{15} is a substituted (C_{6-10})aryl(C_{1-6})alkyl as defined above for each of R^3 and R^4 , respectively;

or the N -oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; and or the pharmaceutically acceptable salts and solvates of such compounds or the N -oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

2. (Currently Amended) The compound of claim 1 in which:

X^2 is $-CHO$, $-C(O)R^5$, $-C(O)CF_3$, $-C(O)CF_2CF_2R^9$, $-CH=CHS(O)_2R^5$,
 $-C(O)CF_2C(O)NR^5R^6$, $-C(O)C(O)NR^5R^6$, $-C(O)C(O)OR^5$, $-C(O)CH_2OR^5$,
 $-C(O)CH_2N(R^6)SO_2R^5$, $-C(O)C(O)N(R^6)(CH_2)_2OR^6$, $-C(O)C(O)N(R^6)(CH_2)_2NR^6$ or
 $-C(O)C(O)R^5$, wherein R^5 is (C_{1-4})alkyl, (C_{6-10})aryl(C_{0-6})alkyl, or (C_{4-10})cycloalkyl(C_{0-6})alkyl, R^6 is hydrogen or (C_{1-6})alkyl and R^9 is halo;

wherein within X^1 any cycloalkyl, or aryl is unsubstituted or substituted with 1 radical selected from $-R^{15}$ and $-X^4C(O)R^{15}$; and wherein X^1 is unsubstituted or substituted further with 1

USAV2001/0081 US NP

PATENT

to 3 radicals independently selected from (C₁₋₆)alkyl, halo-substituted(C₁₋₄)alkyl, -X⁴NR¹²R¹², -X⁴OR¹³ and -X⁴S(O)₂R¹⁴, wherein X⁴, R¹², R¹³, R¹⁴ and R¹⁵ are as defined above;

R¹ and R² are both fluoro; or

R¹ is hydrogen or (C₁₋₆)alkyl and R² is selected from the group consisting of hydrogen, (C₁₋₆)alkyl, -X⁴OR¹³ and -R¹⁵; or R¹ and R² taken together with the carbon atom to which both R¹ and R² are attached form (C₃₋₈)cycloalkylene; wherein R² may be substituted further with (C₁₋₆)alkyl; wherein X⁴, R¹³ and R¹⁵ are as defined above;

R³ and R⁴ are independently is -C(R¹⁶)(R¹⁷)X⁷, wherein R¹⁶ and R¹⁷ are hydrogen, (C₁₋₆)alkyl or fluoro, or R¹⁶ is hydrogen and R¹⁷ is hydroxy and X⁷ is selected from -X⁴SR¹³, -X⁴C(O)R¹³, -X⁴C(O)NR¹²R¹², -R¹⁵, -X⁴OR¹⁵, -X⁴SR¹⁵, -X⁴S(O)₂R¹⁵, -X⁴C(O)R¹⁵ and -X⁴C(O)NR¹⁵R¹², wherein X⁴, R¹², R¹³ and R¹⁵ are as defined above;

wherein within one of R³ or R⁴ any cycloalkyl, heterocycloalkyl, aryl or heteroaryl may be substituted with 1 radical selected from R¹⁵, -X⁴OR¹⁵, -X⁴SR¹⁵, -X⁴S(O)R¹⁵, -X⁴S(O)₂R¹⁵, -X⁴C(O)R¹⁵, -X⁴C(O)OR¹⁵, -X⁴OC(O)R¹⁵, -X⁴NR¹⁵R¹², -X⁴NR¹²C(O)R¹⁵, -X⁴NR¹²C(O)OR¹⁵, -X⁴C(O)NR¹²R¹⁵, -X⁴S(O)₂NR¹⁵R¹², -X⁴NR¹²S(O)₂R¹⁵, -X⁴NR¹²C(O)NR¹⁵R¹² and -X⁴NR¹²C(O)NR¹⁵R¹², wherein X⁴, R¹² and R¹⁵ are as defined above; and wherein each of R³ and R⁴ may be substituted further with 1 to 5 radicals independently selected from (C₁₋₆)alkyl, cyano, halo, halo-substituted(C₁₋₄)alkyl, nitro, -X⁴NR¹²R¹², -X⁴NR¹²C(O)R¹², -X⁴NR¹²C(O)OR¹², -X⁴NR¹²C(O)NR¹²R¹², -X⁴NR¹²C(NR¹²)NR¹²R¹², -X⁴OR¹³, -X⁴SR¹³, -X⁴C(O)OR¹³, -X⁴C(O)R¹³, -X⁴OC(O)R¹³, -X⁴C(O)NR¹²R¹², -X⁴S(O)₂NR¹²R¹², -X⁴NR¹²S(O)₂R¹², -X⁴P(O)(OR¹²)OR¹², -X⁴OP(O)(OR¹²)OR¹², -X⁴S(O)₂R¹⁴ and -X⁴S(O)₂R¹⁴, wherein X⁴, R¹², R¹³ and R¹⁴ are as defined above;

wherein within one of R³ and R⁴ any cycloalkyl, or aryl is unsubstituted or substituted with 1 radical selected from R¹⁵ and -X⁴OR¹⁵; and wherein each of R³ or R⁴ is unsubstituted or substituted further by 1-5 radicals independently selected from (C₁₋₆)alkyl, cyano, halo, halo-substituted(C₁₋₄)alkyl, -X⁴NR¹²C(O)OR¹², -X⁴OR¹³, -X⁴C(O)OR¹², -X⁴C(O)R¹³, -X⁴C(O)NR¹²R¹³, -X⁴NR¹²S(O)₂R¹³ and -X⁴S(O)₂R¹⁴, wherein X⁴, R¹², R¹³, R¹⁴ and R¹⁵ are as defined above;

or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual

USA V2001/0081 US NP

PATENT

isomers and mixtures of isomers thereof.

3. (Currently Amended) A compound of claim 2 in which R^3 and R^4 are independently ~~is~~ $-CH_2X^7$, wherein X^7 is selected from X^4SR^{13} , $-X^4C(O)R^{13}$, $-X^4C(O)NR^{12}R^{12}$, $-R^{15}$, $-X^4OR^{15}$, $-X^4SR^{15}$, $-X^4S(O)_2R^{15}$, $-X^4C(O)R^{15}$ and $-X^4C(O)NR^{15}R^{12}$, wherein X^4 is a bond or (C_{1-6}) alkylene, R^{12} at each occurrence independently is hydrogen or (C_{1-6}) alkyl, R^{13} is hydrogen, (C_{1-6}) alkyl or halo-substituted(C_{1-6})alkyl, R^{14} is (C_{1-6}) alkyl or halo-substituted(C_{1-6})alkyl and R^{15} ~~for R^3 is~~ (C_{3-10}) cycloalkyl(C_{0-6})alkyl, (C_{3-10}) cycloalkyl(C_{0-6})alkyl, morpholinyl, (C_{6-10}) aryl(C_{0-6})alkyl, or (C_{9-12}) bicycloaryl(C_{0-6})alkyl; wherein ~~within~~ R^3 ~~may be substituted and R^4 any cycloalkyl, or~~ ~~aryl or may be substituted with 1 radical selected from R^{15} and X^4OR^{15} , wherein X^4 and R^{15} are as defined above; and wherein R^3 and R^4 may be substituted further by 1 to 5 radicals independently selected from (C_{1-6}) alkyl, cyano, halo, halo substituted($C_{1-4})$ alkyl, $X^4NR^{12}C(O)OR^{12}$, X^4OR^{13} , $X^4C(O)OR^{12}$, $X^4C(O)R^{13}$, $X^4C(O)NR^{12}R^{13}$, $X^4NR^{12}S(O)_2R^{13}$ and $X^4S(O)_2R^{14}$, wherein X^4 , R^{12} , R^{13} and R^{14} are as defined above;~~

or the N -oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N -oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

4. (Currently Amended) A compound of claim 3 in which R^3 is selected from ~~5-bromo-thiophen-2-ylmethyl, 3-cyclohexylpropyl, 2-cyclohexylpropyl, 2-cyclopentylpropyl, 3-phenylpropyl, 3-(2-difluoromethoxy)phenylpropyl, 2-phenylcyclopropylmethyl, 2,2-difluoro-3-phenylpropyl, 1-benzylcyclopropylmethyl, 2-tetrahydro-pyran-4-ylmethyl, 1-isobutylcyclopropylmethyl, thiophen-2-ylmethyl, tetrahydro-pyran-4-ylmethyl, cyclopropylmethylsulfanyl methyl, 2,2-dimethyl-3-phenylpropyl, 4-methyl-[1,2,5]thiadiazol-3-ylmethylsulfonylmethyl, 3-methyl-[1,2,4]thiadiazol-3-ylmethylsulfonylmethyl, thiophen-3-ylmethylsulfonylmethyl, 3-methoxy-5-methyl-isoxazol-4-ylmethylsulfonylmethyl, 2,4-dimethyl-thiazol-5-ylmethylsulfonylmethyl, 2-methyl-oxazol-4-ylmethylsulfonylmethyl, 2-methyl-thiazol-4-ylmethylsulfonylmethyl, 1,2,3-thiadiazol-4-ylmethylsulfonylmethyl, 3-methyl-[1,2,4]thiadiazol-5-ylmethylsulfonylmethyl, 4-methyl-[1,2,5]thiadiazol-3-ylmethylsulfonylmethyl, thiophen-3-ylmethylsulfonylmethyl, tetrahydro-~~

USA2001/0081 US NP

PATENT

pyran-4-ylmethoxy, piperidin-1-ylcarbonyl, thiophene-2-sulfonylmethyl,
3-chloro-2-fluoro-benzylsulfonylmethyl, benzenesulfonylmethyl, benzylsulfonylmethyl,
2-(1,1-difluoro-methoxy)-benzylsulfonylmethyl, 2-benzenesulfonyl-ethyl,
2-(pyridine-2-sulfonyl)-ethyl, 2-(pyridine-4-sulfonyl)-ethyl, 2-benzylsulfonyl-ethyl,
oxy-pyridin-2-ylmethoxy, prop-2-ene-1-sulfonylmethyl,
4-methoxy-benzylsulfonylmethyl, *p*-tolylmethylsulfonylmethyl, 4-chloro-benzylsulfonylmethyl,
o-tolylmethylsulfonylmethyl, 3,5-dimethyl-benzylsulfonylmethyl,
4-trifluoromethyl-benzylsulfonylmethyl, 4-trifluoromethoxy-benzylsulfonylmethyl,
2-bromo-benzylsulfonylmethyl, pyridin-2-ylmethoxy, pyridin-4-ylmethoxy,
naphthalen-2-ylmethoxy, 3-methyl-benzylsulfonylmethyl,
3-trifluoromethyl-benzylsulfonylmethyl, 3-trifluoromethoxy-benzylsulfonylmethyl,
4-fluoro-2-trifluoromethoxy-benzylsulfonylmethyl,
2-fluoro-6-trifluoromethyl-benzylsulfonylmethyl, 3-chloro-benzylsulfonylmethyl,
2-fluoro-benzylsulfonylmethyl, 2-trifluoro-benzylsulfonylmethyl,
2-cyano-benzylsulfonylmethyl, 4-*tert*-butyl-benzylsulfonylmethyl,
2-fluoro-3-methyl-benzylsulfonylmethyl, 3-fluoro-benzylsulfonylmethyl,
4-fluoro-benzylsulfonylmethyl, 2-chloro-benzylsulfonylmethyl,
2,5-difluoro-benzylsulfonylmethyl, 2,6-difluoro-benzylsulfonylmethyl,
2,5-dichloro-benzylsulfonylmethyl, 3,4-dichloro-benzylsulfonylmethyl,
2-(1,1-difluoro-methoxy)-benzylsulfonylmethyl, 2-cyano-benzylsulfonylmethyl,
3-cyano-benzylsulfonylmethyl, 2-trifluoromethoxy-benzylsulfonylmethyl,
2,3-difluoro-benzylsulfonylmethyl, 2,5-difluoro-benzylsulfonylmethyl,
biphenyl-2-ylmethoxy, cyclohexylmethyl, 3-fluoro-benzylsulfonylmethyl,
3,4-difluoro-benzylsulfonylmethyl, 2,4-difluoro-benzylsulfonylmethyl,
2,4,6-trifluoro-benzylsulfonylmethyl, 2,4,5-trifluoro-benzylsulfonylmethyl,
2,3,4-trifluoro-benzylsulfonylmethyl, 2,3,5-trifluoro-benzylsulfonylmethyl,
2,5,6-trifluoro-benzylsulfonylmethyl, 2-chloro-5-trifluoromethylbenzylsulfonylmethyl,
2-methyl-propane-1-sulfonyl, 2-fluoro-3-trifluoromethylbenzylsulfonylmethyl,
2-fluoro-4-trifluoromethylbenzylsulfonylmethyl,
2-fluoro-5-trifluoromethylbenzylsulfonylmethyl,

USA V2001/0081 US NP

PATENT

4-fluoro-3-trifluoromethylbenzylsulfonylmethyl, 2-methoxy-benzylsulfonylmethyl, 3,5-bis-trifluoromethyl-benzylsulfonylmethyl, 4-difluoromethoxy-benzylsulfonylmethyl, 2-difluoromethoxy-benzylsulfonylmethyl, 3-difluoromethoxy-benzylsulfonylmethyl, 2,6-dichloro-benzylsulfonylmethyl, biphenyl-4-ylmethylsulfonylmethyl, 3,5-dimethyl-isoxazol-4-ylmethylsulfonylmethyl, 5-chloro-thiophen-2-ylmethylsulfonylmethyl, 2-[4-(1,1-Difluoro-methoxy)-benzenesulfonyl]-ethyl, 2-[2-(1,1-Difluoro-methoxy)-benzenesulfonyl]-ethyl, 2-[3-(1,1-Difluoro-methoxy)-benzenesulfonyl]-ethyl, 2-(4-trifluoromethoxy-benzenesulfonyl)-ethyl, 2-(3-trifluoromethoxy-benzenesulfonyl)-ethyl, 2-(2-trifluoromethoxy-benzenesulfonyl)-ethyl, (cyanomethyl-methyl-carbamoyl)-methyl, biphenyl-3-ylmethyl, 2-~~oxo~~-2-pyrrolidin-1-yl-ethyl, 2-benzenesulfonyl-ethyl, isobutylsulfanyl-methyl, 2-phenylsulfanyl-ethyl, cyclohexylmethylsulfonylmethyl, 2-cyclohexyl-ethanesulfonyl, benzyl, naphthalen-2-yl, benzylsulfanyl-methyl, 2-trifluoromethyl-benzylsulfanyl-methyl, phenylsulfanyl-ethyl and cyclopropylmethylsulfonylmethyl;

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

5. (Currently Amended) A compound of claim 4 in which R⁴ is selected from 2-trifluorobenzylsulfonylmethyl, 3-phenylsulfanylpropyl, 4-chlorobenzylsulfonylmethyl, thiophen-2-ylsulfonylmethyl, benzylsulfonylmethyl, 4-methylbenzylsulfonylmethyl, 2-phenylsulfonyl-ethyl, 2-pyridin-2-ylsulfonyl-ethyl, 2-pyridin-4-ylsulfonyl-ethyl, 2-benzylsulfonyl-ethyl, 2-(3-difluoromethoxyphenylsulfonyl)-ethyl, naphthalen-2-ylmethylsulfonylmethyl, pyridin-2-ylmethylsulfonylmethyl, 3-methylbenzylsulfonylmethyl, 3-trifluoromethylbenzylsulfonylmethyl, 3-difluoromethoxybenzylsulfonylmethyl, 3-chlorobenzylsulfonylmethyl, 3-fluorobenzylsulfonylmethyl, 4-fluorobenzylsulfonylmethyl, 3-cyanobenzylsulfonylmethyl, 4-cyanobenzylsulfonylmethyl, 3,4-difluorobenzylsulfonylmethyl, benzylsulfonylmethyl, *N*-cyanomethyl-*N*-methylcarbamoylmethyl, 3-bromoethyl, 4-phenylbutyl, 2,2-difluoro-

USA V2001/0081 US NP

PATENT

~~3-phenylpropyl, 4' methylsulfonylaminobiphenyl 3-ylmethyl, 4' ethoxycarbonylaminobiphenyl-3-ylmethyl, 4-methylpiperazin-1-ylcarbonylmethyl, 1-fluoro-2-(4-methylpiperazin-1-yl)-2-exoethyl, 1-hydroxy-4-methylpiperazin-1-yl-2-exoethyl, 1-hydroxy-2-morpholin-4-yl-2-oxoethyl, 1-hydroxy-2-exo-2-pyrrolidin-1-yl-ethyl, 1-fluoro-2-exo-2-pyrrolidin-1-yl-ethyl, 1-fluoro-2-isopropylamino-2-exoethyl, 1-hydroxy-2-isopropylamino-2-exoethyl, 1-fluoro-2-exo-2-piperazin-1-ylethyl, thiophen-3-ylmethylsulfonylmethyl, 4-methyl-[1,2,5]thiadiazol-3-ylmethylsulfonylmethyl, 3-methoxy-5-methyl-isoxazol-4-ylmethylsulfonylmethyl, 2,4-dimethyl-thiazol-5-ylmethylsulfonylmethyl, 2-methyl-exazol-4-ylmethylsulfonylmethyl, 2-methyl-thiazol-4-ylmethylsulfonylmethyl, 2-((1,2,3)thiadiazol-4-ylmethylsulfonyl)-ethyl, 2-(3-methyl-[1,2,4]thiadiazol-5-ylmethylsulfonyl)-ethyl, 2-exo-2-phenyl-ethyl, and 2-morpholin-4-yl-2-oxo-ethyl, 2-benzenesulfonyl-ethyl, 2-naphthalen-2-yl-2-exo-ethyl, 2-benzo[1,3]dioxol-5-yl-2-exo-ethyl, 2-benzo[b]thiophen-2-yl-2-exo-ethyl, 2-biphenyl-4-yl-2-exo-ethyl, 4-benzylsulfonylmethyl, 2-(3-trifluoromethoxy-benzenesulfonyl)-ethyl, 2-exo-2-(4-phenoxy-phenyl)-ethyl, 2-(4-hydroxy-phenyl)-2-exo-ethyl, benzylcarbamoyl-methyl, 4-acetyl-piperazine-1-carboxylic acid-ethyl-ester, cyclohexylcarbamoylmethyl, 2-(3-Chloro-benzo[b]thiophen-2-yl)-2-exo-ethyl, benzenesulfonylmethyl, 2-exo-2-thiophen-2-yl-ethyl, 2-exo-2-thiophen-3-yl-ethyl, naphthalene-2-sulfonylmethyl, 2-(5-methyl-thiophen-2-yl)-2-exo-ethyl, 2-(3-chloro-thiophen-2-yl)-2-exo-ethyl, 5-methyl-thiophene-2-sulfonylmethyl, phenylcarbamoylmethyl, (5,6,7,8-tetrahydro-naphthalen-1-ylcarbamoyl)-methyl, (4-carbamoyl-phenylcarbamoyl)-methyl, (3-carbamoyl-phenylcarbamoyl)-methyl, (butyl-methyl-carbamoyl)-methyl, biphenyl-4-ylmethyl, 2-exo-2-p-tolyl-ethyl, 2-(3-fluoro-4-methoxy-phenyl)-2-exo-ethyl, 2-(4-chloro-phenyl)-2-exo-ethyl, 2-(4-methoxy-phenyl)-2-exo-ethyl, 2-exo-2-(4-trifluoromethoxy-phenyl)-ethyl, 2-(3,4-difluoro-phenyl)-2-exo-ethyl, 2-(3,4-dimethoxy-phenyl)-2-exo-ethyl, 2-(4-fluoro-phenyl)-2-exo-ethyl, 5-methyl-2-exo-hexyl, 3,5-dimethyl-benzylsulfonylmethyl, 4-trifluoromethyl-benzylsulfonylmethyl, 4-trifluoromethoxy-benzylsulfonylmethyl, isopropylcarbamoyl-methyl, 4-dimethylcarbamoylmethyl, pyridin-4-ylcarbamoylmethyl, pyridin-4-ylmethylsulfonylmethyl, pyridin-3-ylmethylsulfonylmethyl, 3,4-dichloro-benzylsulfonylmethyl, pyridin-3-ylcarbamoylmethyl,~~

USA2001/0081 US NP

PATENT

~~4 methoxybenzylsulfonylmethyl, 4 chlorobenzylsulfonylmethyl, thiophene 2-sulfonylmethyl, benzylsulfonylmethyl, p-tolylmethysulfonylmethyl, 2-benzenesulfonyl-ethyl, 2-(pyridine 2-sulfonyl)-ethyl, 2-(pyridine 4-sulfonyl)-ethyl, 2-benzylsulfonyl-ethyl, 2-[3-(1,1-difluoro-methoxy)-benzenesulfonyl]-ethyl, naphthalen-2-ylmethylsulfonylmethyl, pyridin-2-ylmethylsulfonylmethyl, m-tolylmethylsulfonylmethyl, 3-trifluoromethyl-benzylsulfonylmethyl, 3-trifluoromethoxy-benzylsulfonylmethyl, 3-chlorobenzylsulfonylmethyl, 3-fluorobenzylsulfonylmethyl, 4-fluorobenzylsulfonylmethyl, 3-cyano-benzylsulfonylmethyl, 4-cyano-benzylsulfonylmethyl, 3,4-difluorobenzylsulfonylmethyl, (cyanomethyl-methyl-carbamoyl)-methyl, 3-bromo-benzyl, 2-oxo-2-pyrrolidin-1-yl-ethyl, 2-(4'-chlorobiphenyl-4-yl)-2-oxo-ethyl, biphenyl-3-ylmethyl, 2-(1,1-difluoro-methoxy)-benzylsulfonylmethyl, 2-(4-methylsulfonyl-amino-phenyl)-2-oxo-ethyl, 2-oxo-2-piperidin-1-yl-ethyl, 2-(4-methylsulfonyl-piperazin-1-yl)-2-oxo-ethyl, 2-trifluoromethyl-benzylsulfonylmethyl, 4-fluoro-3-trifluoromethyl-benzylsulfonylmethyl, 4-carboxy-benzylsulfonylmethyl, 3,5-bis-trifluoromethyl-benzylsulfonylmethyl, 4-(1,1-difluoro-methoxy)-benzylsulfonylmethyl, 3-(1,1-difluoro-methoxy)-benzylsulfonylmethyl, 5-chlorothiophen-2-ylmethylsulfonylmethyl, 2-[4-(1,1-difluoro-methoxy)-benzenesulfonyl]-ethyl, 2-(4-trifluoromethoxy-benzenesulfonyl)-ethyl, 2-phenylsulfanyl-ethyl, benzylsulfanyl-methyl, 2-trifluoromethyl-benzylsulfanyl-methyl, 2-trifluoromethoxy-benzylsulfanyl-methyl, 2-cyclohexyl-ethyl and isobutylsulfanyl-methyl;~~

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

6. (Currently Amended) The compound of claim 5 in which R¹ is hydrogen or (C₁₋₆)alkyl and R² is hydrogen, -X⁴OR¹³, hetero(C₅₋₁₀)aryl(C₀₋₆)alkyl, (C₅₋₁₀)aryl(C₀₋₆)alkyl or (C₁₋₆)alkyl; or R¹ and R² taken together with the carbon atom to which both R¹ and R² are attached form (C₃₋₈)cycloalkylene; wherein the cycloalkylene is optionally substituted with 1 to 3 (C₁₋₆)alkyl radicals;

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers

USA V2001/0081 US NP

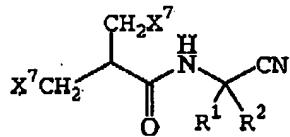
PATENT

and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

7. (Currently Amended) The compound of claim 6 in which R¹ is hydrogen or methyl and R² is methoxymethyl, methoxyethyl, methyl, ethyl, propyl, butyl, or phenethyl, ~~hiophen-2-yl or 5-methyl-furan-2-yl~~; or R¹ and R² taken together with the carbon atom to which both R¹ and R² are attached form cyclopropyl, ~~tetrahydro-pyran-4-yl or 1-methyl-piperidin-4-yl~~;

or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

8. (Previously Presented) The compound of claim 7 of Formula I(a):



I(a)

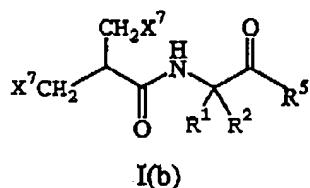
or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

9. (Canceled)

10. (Previously Presented) The compound of claim 7 of Formula I(b):

USA2001/0081 US NP

PATENT



or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

11. (Currently Amended) The compound of claim 10 in which R^5 is 1*H*-benzimidazol-2-yl, benzooxazol-2-yl, oxazolo[4,5-*b*]pyridin-2-yl, benzothiazol-2-yl, 5-phenyl-[1,3,4]oxadiazol-2-yl, 4-(5-pyridin-4-yl-[1,3,4]oxadiazol-2-yl, 5-pyridin-3-yl-[1,3,4]oxadiazol-2-yl, 5-pyridin-3-yl-[1,3,4]oxadiazol-2-yl, pyrimidin-2-yl, ~~pyridazin-3-yl~~, 3-phenyl-[1,2,4]oxadiazol-5-yl, 5-methoxymethyl-[1,3,4]oxadiazol-2-yl, 5-ethyl-[1,3,4]oxadiazol-2-yl, 1,3,4-thiadiazol-2-yl, benzylexycarbonyl, benzylexydicarbonyl, phenyldicarbonyl, 5-methyl-[1,3,4]thiadiazol-2-yl, 5-trifluoromethyl-[1,3,4]oxadiazol-2-yl, 5-methyl-[1,3,4]oxadiazol-2-yl, 5-methyl-[1,2,4]oxadiazol-3-yl, 5-phenyl-[1,2,4]oxadiazol-3-yl, ~~5-thiophen-3-yl-[1,2,4]oxadiazol-3-yl~~, 5-trifluoromethyl-[1,2,4]oxadiazol-3-yl, or 3-methyl-[1,2,4]oxadiazol-5-yl or 3-pyrazin-2-yl;

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

12. (Currently Amended) The compound of claim 11 selected from the group consisting of ~~N[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl]-3-benzylsulfonyl-2-benzylsulfonylmethyl-propionamide~~; ~~N[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl]-3-(2-trifluoromethyl-benzylsulfonyl)-2-(2-trifluoromethyl-benzylsulfonylmethyl)-propionamide~~; ~~N[(S)-1-(1-benzooxazol-2-yl-methanoyl)-pentyl]-4-(2-methoxy-benzenesulfonyl)-2-[2-(2-methoxy-benzenesulfonyl)-ethyl]-butyramide~~; ~~4-Benzenesulfonyl-2-(2-benzenesulfonyl-ethyl)-N[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl]-butyramide~~; ~~(R)-N[(S)-1-(1-benzooxazol-2-yl-~~

USA V2001/0081 US NP

PATENT

~~methanoyl) butyl] 2-cyclohexylmethyl 3-benzylsulfonyl-propionamide; N-[(S)-1-(1-benzothiazol-2-yl-methanoyl)-propyl]-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethylbutyramide; N-[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl] 3-cyclohexyl-2-cyclohexylmethyl-propionamide; N-[(S)-1-(1-Benzooxazol-2-yl-methanoyl)-butyl] 3-isobutylsulfanyl-2-isobutylsulfanyl-methyl-propionamide; N-[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl] 3-benzylsulfanyl-2-benzylsulfanyl-methyl-propionamide; N-[(S)-1-(1-benzooxazol-2-yl-methanoyl)-butyl] 4-phenylsulfanyl-2-(2-phenylsulfanyl-ethyl) butyramide; N-[(S)-1-(1-benzooxazol-2-yl-methanoyl)-propyl]-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethylbutyramide; N-[(S)-1-(1-Benzooxazol-2-yl-methanoyl)-butyl]-2-[2-(1,1-difluoro-methoxy)-benzylsulfonylmethyl]-3-benzylsulfonyl-propionamide; 4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-N-[(S)-1-[1-(3-phenyl-[1,2,4]oxadiazol-5-yl)-methanoyl]-propyl]-butyramide; N-[(S)-1-(1-Benzooxazol-2-yl-methanoyl)-butyl]-4-oxo-2-benzylsulfonylmethyl-butyramide; N-[(S)-1-(1-Benzooxazol-2-yl-methanoyl)-butyl]-2-[2-(1,1-difluoro-methoxy)-benzylsulfonylmethyl]-3-benzylsulfonyl-propionamide; 4-Morpholin-4-yl-4-oxo-N-[1-(2-oxo-2-phenyl-acetyl)-pentyl]-2-benzylsulfonylmethyl-butyramide; N-[(1,1-Dimethyl-2-oxazolo[4,5-b]pyridin-2-yl-2-oxo-ethyl)-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl]-butyramide; N-[1-(5-Ethyl-[1,3,4]oxadiazole-2-carbonyl)-butyl]-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyramide; N-[1-(5-Ethyl-[1,3,4]oxadiazole-2-carbonyl)-butyl]-4-oxo-2-benzylsulfonyl-methyl-4-piperidin-1-yl-butyramide; N-[1-(5-Ethyl-[1,3,4]oxadiazole-2-carbonyl)-butyl]-4-oxo-2-benzylsulfonyl-methyl-4-piperidin-1-yl-butyramide; N-[1-(5-Methoxymethyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyramide; N-[1-(5-Methoxymethyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonylmethyl-4-piperidin-1-yl-butyramide; N-[1-(5-Methoxymethyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonylmethyl-4-piperidin-1-yl-butyramide; 4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-N-[1-(5-phenyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; 4-Oxo-2-benzylsulfonylmethyl-N-[1-(5-phenyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-piperidin-1-yl-butyramide; 4-Oxo-2-benzylsulfonylmethyl-N-[1-(5-phenyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-piperidin-1-yl-butyramide; 4-Morpholin-4-yl-N-[1-(oxazolo[4,5-b]pyridine-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonylmethyl-butyramide; N-[1-(Oxazolo[4,5-b]pyridine-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonyl-methyl-4-piperidin-1-yl-butyramide; N-[1-(Oxazolo[4,5-b]pyridine-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonyl-methyl-4-piperidin-1-yl-butyramide; 4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-N-[1-(5-pyridin-4-yl-~~

USA V2001/0081 US NP

PATENT

[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; 4-Oxo-2-benzylsulfonylmethyl-4-piperidin-1-yl-N-[1-(5-pyridin-4-yl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; 4-Oxo-2-benzylsulfonylmethyl N-[1-(5-pyridin-4-yl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-4-pyrrolidin-1-yl butyramide; 4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-N-[1-(5-pyridin-3-yl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; N-[1-(Benzooxazole-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonylmethyl-4-piperidin-1-yl butyramide; N-[1-(Benzooxazole-2-carbonyl)-propyl]-4-oxo-2-benzylsulfonylmethyl-4-pyrrolidin-1-yl butyramide; N-[1-(Benzooxazole-2-carbonyl)-propyl]-2-cyclohexylmethyl-4-morpholin-4-yl-4-oxo-butyramide; 2-Cyclohexylmethyl-4-morpholin-4-yl-N-[1-(oxazolo[4,5-b]pyridine-2-carbonyl)-propyl]-4-oxo-butyramide; 2-Cyclohexylmethyl-N-[1-(5-ethyl-[1,3,4]oxadiazole-2-carbonyl)-butyl]-4-morpholin-4-yl-4-oxo-butyramide; N-(2-Benzooxazol-2-yl-1-methoxymethyl-2-oxo-ethyl)-2-(2-difluoromethoxy-benzylsulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyramide; N-[1-(Benzooxazole-2-carbonyl)-propyl]-2-(2-cyclohexyl-ethyl)-4-morpholin-4-yl-4-oxo-butyramide; 2-(2-Cyclohexyl-ethyl)-4-morpholin-4-yl-N-[1-(oxazolo[4,5-b]pyridine-2-carbonyl)-propyl]-4-oxo-butyramide; 2-(2-Cyclohexyl-ethyl)-4-morpholin-4-yl-4-oxo-N-[1-(5-phenyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; 2-(2-Difluoromethoxy-benzylsulfonylmethyl)-4-morpholin-4-yl-4-oxo-N-[1-(5-phenyl-[1,3,4]oxadiazole-2-carbonyl)-propyl]-butyramide; 2-(2-Difluoromethoxy-benzylsulfonylmethyl)-N-[1-(5-ethyl-[1,3,4]oxadiazole-2-carbonyl)-butyl]-4-morpholin-4-yl-4-oxo-butyramide; N-[1-(Benzooxazole-2-carbonyl)-propyl]-2-(2-difluoromethoxy-benzyl-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyramide;

~~2-(2-Morpholin-4-yl-2-oxo-ethyl)-5-phenyl-pentanoic acid, 1-(benzooxazole-2-carbonyl)-propyl- amide; (R)-2-Cyclohexylmethyl-4-morpholin-4-yl-4-oxo-N-[(S)-1-(5-phenyl-1,2,4-oxadiazole-3-carbonyl)-propyl]-butyramide; 2-(2-Morpholin-4-yl-2-oxo-ethyl)-5-phenyl-pentanoic acid, (S)-1-(5-phenyl-[1,2,4]oxadiazole-3-carbonyl)-propyl- amide; 4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-N-[(S)-1-(5-phenyl-1,2,4-oxadiazole-3-carbonyl)-propyl]-butyramide; (R)-2-Cyclohexylmethyl-4-morpholin-4-yl-4-oxo-N-[(S)-1-(3-phenyl-1,2,4-oxadiazole-5-carbonyl)-propyl]-butyramide; 4-Morpholin-4-yl-N-[1-(oxazole-2-carbonyl)-3-phenyl-propyl]-4-oxo-2-benzylsulfonylmethyl-butyramide; N-(1,1-Dimethyl-2-oxazol-2-yl-2-oxo-ethyl)-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyramide; N-4-Isopropyl N-1-[exazole-2-carbonyl]-3-phenyl-propyl]-2-benzylsulfonylmethyl-succinamide; 2-(2-~~

USA V2001/0081 US NP

PATENT

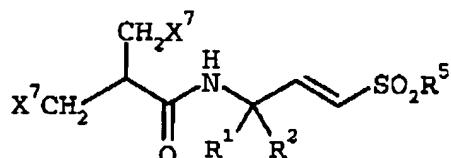
Difluoromethoxy-benzylsulfonylmethyl)-4-morpholin-4-yl-N-[1-(oxazole-2-carbonyl)-3-phenyl-propyl]-4-oxo-butyramide; 2-(2-Methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-N-[1-(oxazole-2-carbonyl)-3-phenyl-propyl]-4-oxo-butyramide; 2-Cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-N-[1-(oxazole-2-carbonyl)-3-phenyl-propyl]-4-oxo-butyramide; *N*-[1-(Benzoxazole-2-carbonyl)-butyl]-2-benzylsulfonyl-3-(tetrahydro-pyran-4-yloxymethyl)-propionamide; *N*-[1-(Benzoxazole-2-carbonyl)-butyl]-3-ethanesulfonyl-2-(tetrahydro-pyran-4-yloxymethyl)-propionamide; *N*-(1-Benzenesulfonyl-3-oxo-azepan-4-yl)-2-cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-4-oxo-butyramide; 2-Cyclopropylmethylsulfonylmethyl-*N*-(S)-1-[*(R)*-hydroxy-(3-phenyl-1,2,4-oxadiazol-5-yl)-methyl]-propyl)-4-morpholin-4-yl-4-oxo-butyramide; *N*-(S)-1-[*(R)*-hydroxy-(3-phenyl-1,2,4-oxadiazol-5-yl)-methyl]-propyl)-2-(2-methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyramide; 2-(2-Morpholin-4-yl-2-oxo-ethyl)-5-phenyl-pentanoic acid (*(S)* 1-[*(R)*-hydroxy-(3-phenyl-1,2,4-oxadiazol-5-yl)-methyl]-propyl)-amide; 2-Cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-4-oxo-*N*-(S)-1-(3-phenyl-1,2,4-oxadiazole-5-carbonyl)-propyl]-butyramide; 2-(2-methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-*N*-(S)-1-(3-phenyl-1,2,4-oxadiazole-5-carbonyl)-propyl]-butyramide; 2-(2-Morpholin-4-yl-2-oxo-ethyl)-5-phenyl-pentanoic acid, (*S*) 1-(3-phenyl-1,2,4-oxadiazole-5-carbonyl)-propyl]-amide; *N*-[(1*S*)-1-(Benzoxazol-2-yl-hydroxy-methyl)-3-phenyl-propyl]-2-cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-4-oxo-butyramide; (*R*) 2-((*S*)-1-Hydroxy-2-morpholin-4-yl-2-oxo-ethyl)-5-phenyl-pentanoic acid, 1-(benzoxazole-2-carbonyl)-propyl]-amide; (*R*) 5-(2-Difluoromethoxy-phenyl)-2-((*S*)-1-hydroxy-2-morpholin-4-yl-2-oxo-ethyl)-pentanoic acid, 1-(benzoxazole-2-carbonyl)-propyl]-amide; and 4-Morpholin-4-yl-*N*-[1-(oxazole-2-carbonyl)-cyclopropyl]-4-oxo-2-benzylsulfonylmethyl-butyramide;

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

13. (Previously Presented) The compound of claim 7 of Formula I(c):

USA V2001/0081 US NP

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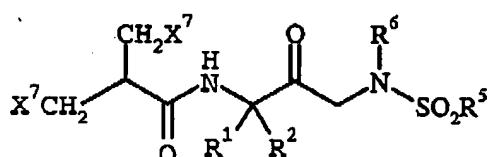
I(c)

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

14. (Previously Presented) The compound of claim 13 in which R^5 is phenyl; or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

15. (Canceled)

16. (Previously Presented) The compound of claim 7 of Formula I(d):



I(d)

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

USA2001/0081 US NP

PATENT

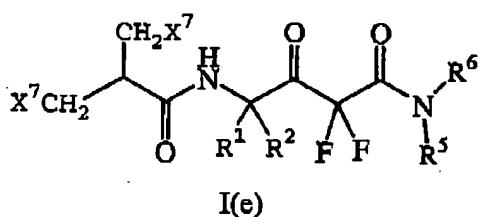
17. (Previously Presented) The compound of claim 16 in which R⁵ is phenyl and R⁶ is hydrogen;

or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

18. (Previously Presented) The compound of claim 17 namely *N*-(3-benzenesulfonylamino-2-oxo-propyl)-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butamide;

or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

19. (Previously Presented) The compound of claim 7 of Formula I(e):



or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

20. (Previously Presented) The compound of claim 19 in which R⁵ and R⁶ is methyl;

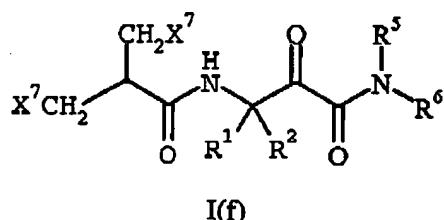
or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

USA2001/0081 US NP

PATENT

21. (Canceled).

22. (Previously Presented) The compound of claim 7 of Formula I(f):



or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

23. (Currently Amended) The compound of claim 22 in which R^5 is methyl, benzyl, phenethyl, cyclohexyl, methoxyethyl, dimethylaminoethyl, tetrahydro-pyran-4-yl, ~~± methylsulfonyl piperidin-4-yl, 4-methyl piperazin-1-yl, morpholin-4-ylethyl, pyridin-2-yl, pyridin-2-ylmethyl or oxazol-2-ylmethyl; or R^6 is hydrogen or methyl; or R^5 and R^6 together with the nitrogen atom to which both R^5 and R^6 are attached form morpholine-4-yl, pyrrolidin-1-yl, 4-dimethylamino piperazin-1-yl, 4-hydroxy piperazin-1-yl, 4-pyridin-2-yl piperazin-1-yl, 4-benzoyl piperazin-1-yl or 3-oxo piperazin-1-yl;~~

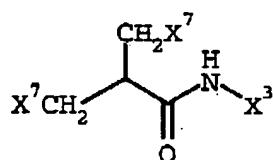
or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

24. (Canceled)

25. (Previously Presented) The compound of claim 7 of Formula I(g):

USA V2001/0081 US NP

PATENT



I(g)

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such compounds or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

26. (Cancelled)

27. (Previously Presented) The compound of claim 23 selected from the group consisting of 3-Hydroxy-4-(4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyrylamino)-azepane-1-carboxylic acid tert-butyl ester; 4-(2-Cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-4-oxo-butyrylamino)-3-hydroxy-azepane-1-carboxylic acid tert-butyl ester; 3-Hydroxy-4-[2-(2-methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyrylamino]-azepane-1-carboxylic acid tert-butyl ester; 4-(4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyrylamino)-3-oxo-azepane-1-carboxylic acid tert-butyl ester; 4-(2-Cyclopropylmethylsulfonylmethyl-4-morpholin-4-yl-4-oxo-butyrylamino)-3-oxo-azepane-1-carboxylic acid tert-butyl ester; 4-[2-(2-Methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyrylamino]-3-oxo-azepane-1-carboxylic acid tert-butyl ester; *N*-(1-Benzenesulfonyl-3-oxo-azepan-4-yl)-4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyramide; *N*-(1-Benzenesulfonyl-3-oxo-azepan-4-yl)-2-(2-methyl-propane-1-sulfonylmethyl)-4-morpholin-4-yl-4-oxo-butyramide; 3-(4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyrylamino)-4-oxo-pyrrolidine-1-carboxylic acid tert-butyl ester; 4-(4-Morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butyrylamino)-3-oxo-azepane-1-carboxylic acid benzyl ester; and acetic acid (2*S*,3*S*)-3-(4-morpholin-4-yl-4-oxo-2-benzylsulfonylmethyl-butanoylamino)-4-oxo-azetidin-2-yl ester;

or the *N*-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof; or the pharmaceutically acceptable salts and solvates of such

USA V2001/0081 US NP

PATENT

compounds or the N-oxide derivatives, prodrug derivatives, protected derivatives, individual isomers and mixtures of isomers thereof.

28. (Previously Presented) A pharmaceutical composition comprising a therapeutically effective amount of a compound of claim 1 in combination with a pharmaceutically acceptable excipient.

29-31. (Canceled)